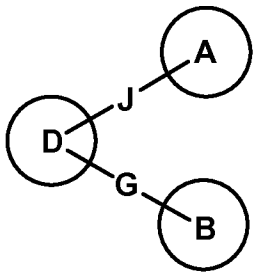
AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions and listings of claims in the application:

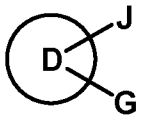
LISTING OF CLAIMS:

1. (currently amended): A compound of formula (I):

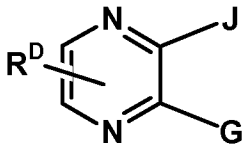


(I)

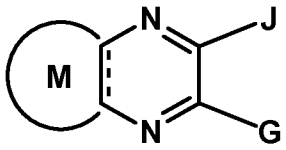
wherein



is



or



wherein R^D represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) halogen, (6) cyano, (7) nitro, (8) $-CONR^7R^8$, (9) $-COOR^9$, (10) Cyc1 or (11) C_{1-8} alkyl substituted with 1 to 5 groups selected from (a) $-CONR^7R^8$, (b) $-COOR^9$, (c) $-OR^{10}$, (d) $-NR^{11}R^{12}$, (e) halogen, and (f) Cyc1;

R^7 and R^8 each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc2, (6) $-OR^{13}$ or (7) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl

substituted with 1 to 5 groups selected from (a) $-OR^{13}$, (b) $-NR^{14}R^{15}$, (c) $-NR^{16}COR^{17}$, (d) halogen, (e) CF_3, and (f) Cyc2; or

R^7 and R^8 are taken together with the adjacent nitrogen atom to represent a pyrrolidine ring, a piperidine ring, an azepane ring, a piperazine ring, a morpholine ring, a thiomorpholine ring, a 2,5-dihydropyrrole ring or a 1,2,3,6-tetrahydropyridine ring which may be substituted with (a) C_{1-8} alkyl, (b) halogen, (c) hydroxyl, or (d) C_{1-8} alkyl substituted with hydroxyl;

R^{13} to R^{17} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc1, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc1;

R^9 to R^{12} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc1, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc1;

Cyc1 represents a benzene ring, naphthalene ring, thiophen ring, 1,3-benzodioxole ring or phenanthrene ring, wherein ~~Cycl1~~ Cyc1 may be substituted with 1 to 5 of R^{18};

R^{18} represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{19}$, (10) $-SR^{20}$, (11) $-NR^{21}R^{22}$, (12) $-COR^{23}$, (13) $-COOR^{24}$, (14) $-NR^{25}COR^{26}$, (15) $-CONR^{27}R^{28}$, (16) Cyc2, or (17) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) halogen, (b) cyano, (c) nitro, (d) trifluoromethyl, (e) trifluoromethoxy, (f) $-OR^{19}$, (g) $-SR^{20}$, (h) $-NR^{21}R^{22}$, (i) $-COR^{23}$, (j) $-COOR^{24}$, (k) $-NR^{25}COR^{26}$, (l) $-CONR^{27}R^{28}$, and (m) Cyc2;

R^{19} to R^{28} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc2, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc2;

Cyc2 represents a benzene, cyclobutane, cyclopentane, cyclohexane, pyridine, pyrimidine, tetrahydrofuran, thiazole, oxazole, isoxazole, pyrazole or 1,2,4-thiadiazole, wherein Cyc2 may be substituted with 1 to 5 of R^{29};

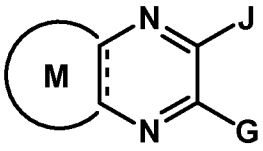
R^{29} represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) hydroxyl, (8) trifluoromethyl, (9) trifluoromethoxy, or (10) $-OR^{100}$;

R^{100} represents C_{1-8} alkyl;

---- represents a single bond or a double bond; and

M represents a 5- to 6-membered ~~3- to 11-membered monocyclic or bicyclic~~ cyclic group which may be substituted;

wherein



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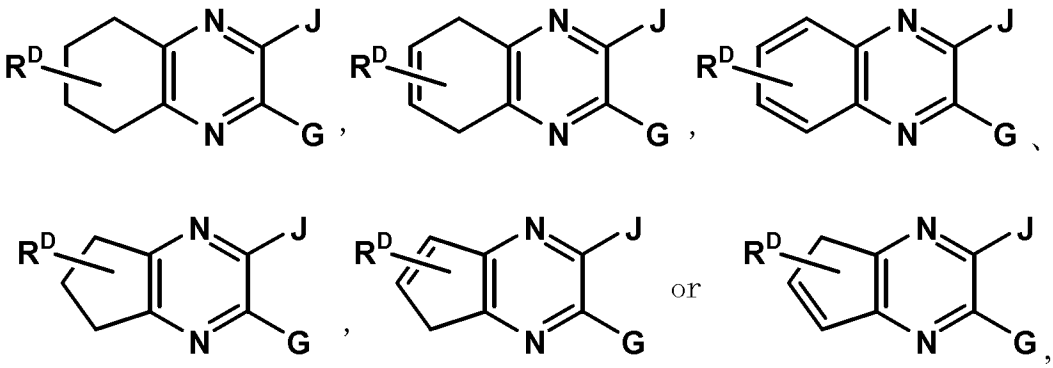


ring A is a benzene ring, a naphthalene ring, a pyridine ring, a pyrazole ring, a dioxaindane ring, a benzodioxane ring, a cyclopropane ring, a cyclopentane ring, a furan ring, a thiophene ring, a tetrahydrofuran ring, a piperidine ring or a morpholine ring which may be substituted with (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{31}$, (10) $-NR^{32}R^{33}$, (11) $-NR^{34}COR^{35}$, (12) Cyc3, or (13) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) halogen, (b) cyano, (c) nitro, (d) trifluoromethyl, (e) trifluoromethoxy, (f) $-OR^{31}$, (g) $-NR^{32}COR^{33}$, (h) $-NR^{34}COR^{35}$, and (i) Cyc3;

R^{31} to R^{35} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc3, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) Cyc3, (b) $-OR^{36}$ and (c) $-NR^{37}R^{38}$;

R^{36} to R^{38} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) $-OR^{39}$, or (4) $-NR^{40}R^{41}$;

R^{39} to R^{41} each independently represents hydrogen or C_{1-8} alkyl;

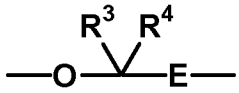
Cyc3 represents a benzene ring, a piperidine ring or a morpholine ring;

ring B is a benzene ring, a pyridine ring, a thiophene ring, ~~a naphthalene ring,~~ a pyrrole ring, a pyrazole ring, an isoxazole ring, a thiazole ring, ~~a benzothiophene ring,~~ an imidazole ring or a furan ring which may be substituted with (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{42}$, (10) $-NR^{43}R^{44}$, (11) $-SR^{101}$, (12) $-SO_2R^{102}$, (13) $-COR^{103}$, (14) $-COOR^{104}$, (15) Cyc2, or (16) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) $-COOR^{104}$, (b) $-NR^{105}COR^{106}$, and (c) Cyc2;

R^{42} to R^{44} and R^{101} to R^{106} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) Cyc2, or (4) $-COR^{107}$, or (5) C_{1-8} alkyl substituted with 1 to 5 halogen atoms;

R^{107} represents C_{1-8} alkyl;

J is



wherein R^3 and R^4 each independently represents hydrogen or C_{1-8} alkyl; and

E represents a bond or a spacer having 1 to 6 atoms in its main chain; and

G is $-NR^{T1}-SO_2-$

wherein R^{T1} represents hydrogen, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl or a 3- to 8-membered cyclic group;

or a salt thereof.

Claims 2-24. (canceled).

25. (previously presented): The compound according to claim 1, wherein R^3 and R^4 each independently represents hydrogen or methyl.

26. (previously presented): The compound according to claim 1, wherein E is a bond.

27. (previously presented): The compound according to claim 1, wherein E is a spacer having 1 to 6 atoms in its main chain.

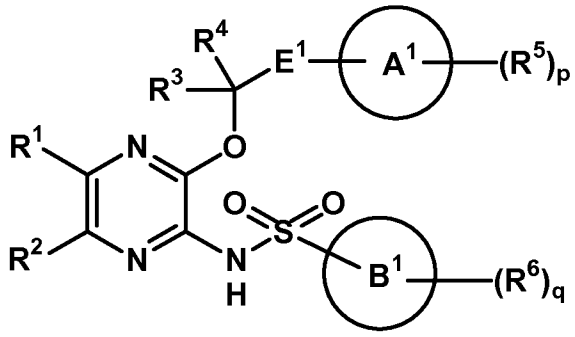
28. (original): The compound according to claim 27, wherein E is C_{1-4} alkylene or C_{1-3} alkyleneoxy.

29. (original): The compound according to claim 28, wherein E is methylene or methylenoxy.

Claims 30-31. (canceled).

32. (previously presented): The compound according to claim 1, wherein G is -NH-SO_2-.

33. (currently amended): The compound according to claim 1, wherein the compound is a compound of formula (A):



(A)

wherein R^1 and R^2 each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) halogen, (6) cyano, (7) nitro, (8) -$CONR^7R^8$, (9) -$COOR^9$, (10) Cyc1 or (11) C_{1-8} alkyl substituted with 1 to 5 groups selected from (a) -$CONR^7R^8$, (b) -$COOR^9$, (c) -OR^{10}, (d) -$NR^{11}R^{12}$, (e) halogen, and (f) Cyc1; or

R^1 and R^2 are taken together to represent C_{3-4} alkylene, -CH=CH-CH_2-, -CH_2-CH=CH-, -CH=CH-CH=CH- or -CH=CH-CH_2-CH_2-, wherein the carbocyclic ring to be formed may be substituted with C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl, C_{1-8} alkoxy, halogen, cyano, nitro or hydroxyl, wherein R^7 and R^8 each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc2, (6) -OR^{13} or (7) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) -OR^{13}, (b) -$NR^{14}R^{15}$, (c) -$NR^{16}COR^{17}$, (d) halogen, (e) CF_3, and (f) Cyc2; or

R^7 and R^8 are taken together with the adjacent nitrogen atom to represent a pyrrolidine ring, a piperidine ring, an azepane ring, a piperazine ring, a morpholine ring, a thiomorpholine ring, a 2,5-dihydropyrrole ring or a 1,2,3,6-tetrahydropyridine ring which may be substituted with (a) C_{1-8} alkyl, (b) halogen, (c) hydroxyl, or (d) C_{1-8} alkyl substituted with hydroxyl;

R^{13} to R^{17} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc1, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc1;

R^9 to R^{12} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc1, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc1;

Cyc1 represents a benzene ring, a naphthalene ring, thiophene ring, 1,3-benzodioxole ring or phenanthrene ring, wherein Cyc1 may be substituted with 1 to 5 of R^{18};

R^{18} represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) -OR^{19}, (10) -SR^{20}, (11) -$NR^{21}R^{22}$, (12) -COR^{23}, (13) -$COOR^{24}$, (14) -$NR^{25}COR^{26}$, (15) -$CONR^{27}R^{28}$, (16) Cyc2, or (17) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) halogen, (b) cyano, (c) nitro, (d) trifluoromethyl, (e) trifluoromethoxy, (f) -OR^{19}, (g) -SR^{20}, (h) -$NR^{21}R^{22}$, (i) -COR^{23}, (j) -$COOR^{24}$, (k) -$NR^{25}COR^{26}$, (l) -$CONR^{27}R^{28}$, and (m) Cyc2;

R^{19} to R^{28} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc2, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc2;

Cyc2 represents a benzene, cyclobutane, cyclopentane, cyclohexane, pyridine, pyrimidine, tetrahydrofuran, thiazole, oxazole, isoxazole, pyrazole or 1,2,4-thiadiazole, wherein Cyc2 may be substituted with 1 to 5 of R^{29};

R^{29} represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) hydroxyl, (8) trifluoromethyl, (9) trifluoromethoxy, or (10) $-OR^{100}$;

R^{100} represents C_{1-8} alkyl.;

R^3 and R^4 each independently represents hydrogen or C_{1-8} alkyl;

E^1 represents a bond or C_{1-6} alkylene, wherein a carbon atom in the alkylene group may be substituted with oxygen, sulfur, or $-NR^{30}-$;

R^{30} represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) phenyl, or (5) C_{1-8} alkyl substituted with phenyl;

ring A^1 is a benzene ring, a naphthalene ring, a pyridine ring, a pyrazole ring, a dioxaindane ring, a benzodioxane ring, a cyclopropane ring, a cyclopentane ring, a furan ring, a thiophene ring, a tetrahydrofuran ring, a piperidine ring or a morpholine ring;

R^5 represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{31}$, (10) $-NR^{32}R^{33}$, (11) $-NR^{34}COR^{35}$, (12) Cyc3, or (13) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) halogen, (b) cyano, (c) nitro, (d) trifluoromethyl, (e) trifluoromethoxy, (f) $-OR^{31}$, (g) $-NR^{32}COR^{33}$, (h) $-NR^{34}COR^{35}$, and (i) Cyc3;

R^{31} to R^{35} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc3, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) Cyc3, (b) $-OR^{36}$ and (c) $-NR^{37}R^{38}$;

R^{36} to R^{38} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) $-OR^{39}$, or (4) $-NR^{40}R^{41}$;

R^{39} to R^{41} each independently represents hydrogen or C_{1-8} alkyl;

Cyc3 represents a benzene ring, a piperidine ring or a morpholine ring;

ring B^1 is a benzene ring, a pyridine ring, a thiophene ring, ~~a naphthalene ring,~~ a pyrrole ring, a pyrazole ring, an isoxazole ring, a thiazole ring, ~~a benzothiophene ring,~~ an imidazole ring or a furan ring;

R^6 represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{42}$, (10) $-NR^{43}R^{44}$, (11) $-SR^{101}$, (12) $-SO_2R^{102}$, (13) $-COR^{103}$, (14) $-COOR^{104}$, (15) Cyc2, or (16) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) $-COOR^{104}$, (b) $-NR^{105}COR^{106}$, and (c) Cyc2;

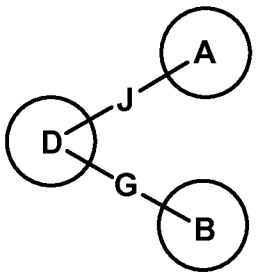
R^{42} to R^{44} and R^{101} to R^{106} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) Cyc2, or (4) $-COR^{107}$, or (5) C_{1-8} alkyl substituted with 1 to 5 halogen atoms;

R^{107} represents C_{1-8} alkyl; and

p and q each independently represents 0 or an integer of 1 to 5.

34. (withdrawn): A prodrug for the compound according to claim 1.

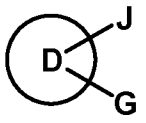
35. (currently amended): A pharmaceutical composition which comprises the compound of formula (I):



(I)

;

wherein



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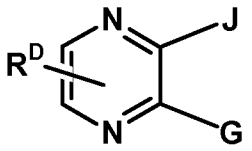
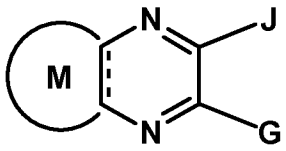


or



wherein R^D represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) halogen, (6) cyano, (7) nitro, (8) $-CONR^7R^8$, (9) $-COOR^9$, (10) Cyc1 or (11) C_{1-8} alkyl substituted with 1 to 5 groups selected from (a) $-CONR^7R^8$, (b) $-COOR^9$, (c) $-OR^{10}$, (d) $-NR^{11}R^{12}$, (e) halogen, and (f) Cyc1;

R^7 and R^8 each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc2, (6) $-OR^{13}$ or (7) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) $-OR^{13}$, (b) $-NR^{14}R^{15}$, (c) $-NR^{16}COR^{17}$, (d) halogen, (e) CF_3, and (f) Cyc2; or

R^7 and R^8 are taken together with the adjacent nitrogen atom to represent a pyrrolidine ring, a piperidine ring, an azepane ring, a piperazine ring, a morpholine ring, a thiomorpholine ring, a 2,5-dihydropyrrole ring or a 1,2,3,6-tetrahydropyridine ring which may be substituted with (a) C_{1-8} alkyl, (b) halogen, (c) hydroxyl, or (d) C_{1-8} alkyl substituted with hydroxyl;

R^{13} to R^{17} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc1, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc1;

R^{9} to R^{12} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc1, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc1;

Cyc1 represents a benzene ring, naphthalene ring, thiophene ring, 1,3-benzodioxole ring or phenanthrene ring, wherein Cyc1 may be substituted with 1 to 5 of R^{18};

R^{18} represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{19}$, (10) $-SR^{20}$, (11) $-NR^{21}R^{22}$, (12) $-COR^{23}$, (13) $-COOR^{24}$, (14) $-NR^{25}COR^{26}$, (15) $-CONR^{27}R^{28}$, (16) Cyc2, or (17) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) halogen, (b) cyano, (c) nitro, (d) trifluoromethyl, (e) trifluoromethoxy, (f) $-OR^{19}$, (g) $-SR^{20}$, (h) $-NR^{21}R^{22}$, (i) $-COR^{23}$, (j) $-COOR^{24}$, (k) $-NR^{25}COR^{26}$, (l) $-CONR^{27}R^{28}$, and (m) Cyc2;

R^{19} to R^{28} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc2, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with Cyc2;

Cyc2 represents a benzene, cyclobutane, cyclopentane, cyclohexane, pyridine, pyrimidine, tetrahydrofuran, thiazole, oxazole, isoxazole, pyrazole or 1,2,4-thiadiazole, wherein Cyc2 may be substituted with 1 to 5 of R^{29};

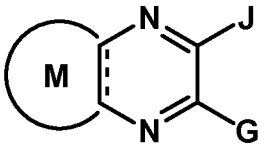
R^{29} represents (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) hydroxyl, (8) trifluoromethyl, (9) trifluoromethoxy, or (10) $-OR^{100}$;

R^{100} represents C_{1-8} alkyl;

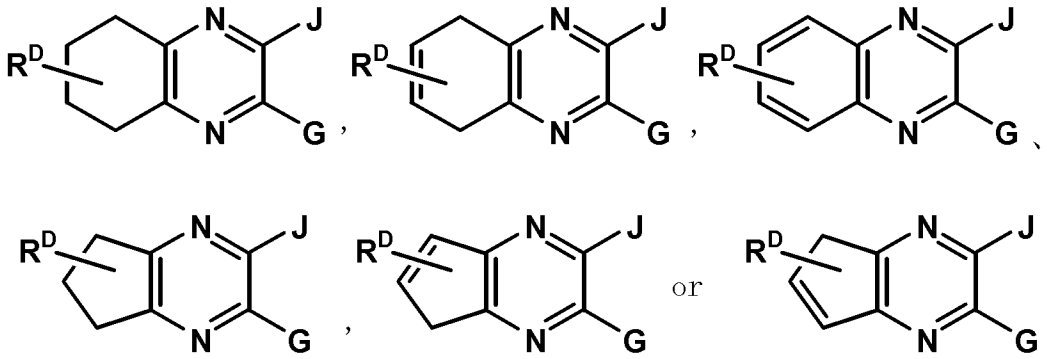
---- represents a single bond or a double bond; and

M represents a 5- to 6-membered~~3- to 11-membered monocyclic or bicyclic~~ cyclic group which may be substituted;

wherein



is



ring A is a benzene ring, a naphthalene ring, a pyridine ring, a pyrazole ring, a dioxaindane ring, a benzodioxane ring, a cyclopropane ring, a cyclopentane ring, a furan ring, a thiophene ring, a tetrahydrofuran ring, a piperidine ring or a morpholine ring which may be substituted with (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{31}$, (10) $-NR^{32}R^{33}$, (11) $-NR^{34}COR^{35}$, (12) Cyc3, or (13) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) halogen, (b) cyano, (c) nitro, (d) trifluoromethyl, (e) trifluoromethoxy, (f) $-OR^{31}$, (g) $-NR^{32}COR^{33}$, (h) $-NR^{34}COR^{35}$, and (i) Cyc3;

R^{31} to R^{35} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) C_{2-8} alkenyl, (4) C_{2-8} alkynyl, (5) Cyc3, or (6) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) Cyc3, (b) $-OR^{36}$ and (c) $-NR^{37}R^{38}$;

R^{36} to R^{38} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) $-OR^{39}$, or (4) $-NR^{40}R^{41}$;

R^{39} to R^{41} each independently represents hydrogen or C_{1-8} alkyl;

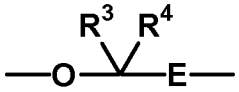
Cyc3 represents a benzene ring, a piperidine ring or a morpholine ring;

ring B is a benzene ring, a pyridine ring, a thiophene ring, ~~a naphthalene ring,~~ a pyrrole ring, a pyrazole ring, an isoxazole ring, a thiazole ring, ~~a benzothiophene ring,~~ an imidazole ring or a furan ring which may be substituted with (1) C_{1-8} alkyl, (2) C_{2-8} alkenyl, (3) C_{2-8} alkynyl, (4) halogen, (5) cyano, (6) nitro, (7) trifluoromethyl, (8) trifluoromethoxy, (9) $-OR^{42}$, (10) $-NR^{43}R^{44}$, (11) $-SR^{101}$, (12) $-SO_2R^{102}$, (13) $-COR^{103}$, (14) $-COOR^{104}$, (15) Cyc2, or (16) C_{1-8} alkyl, C_{2-8} alkenyl or C_{2-8} alkynyl substituted with 1 to 5 groups selected from (a) $-COOR^{104}$, (b) $-NR^{105}COR^{106}$, and (c) Cyc2;

R^{42} to R^{44} and R^{101} to R^{106} each independently represents (1) hydrogen, (2) C_{1-8} alkyl, (3) Cyc2, or (4) $-COR^{107}$, or (5) C_{1-8} alkyl substituted with 1 to 5 halogen atoms;

R^{107} represents C_{1-8} alkyl;

J is



wherein R^3 and R^4 each independently represents hydrogen or C_{1-8} alkyl; and

E represents a bond or a spacer having 1 to 6 atoms in its main chain;

G is $-NR^{T1}-SO_2-$

wherein R^{T1} represents hydrogen, C_{1-8} alkyl, C_{2-8} alkenyl, C_{2-8} alkynyl or a 3- to 8-membered cyclic group;

or a salt thereof and a pharmaceutically acceptable carrier.

Claims 36-49. (canceled).